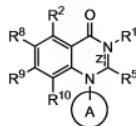


Amendments to the Claims

1. (Canceled) A compound of the structure:



or a pharmaceutically acceptable salt thereof, wherein

z is a single or double bond;

A is

a) an aryl ring, wherein any stable aryl ring atom is independently unsubstituted or substituted with

- 1) halogen,
- 2) NO₂,
- 3) CN,
- 4) CR⁴⁶=C(R⁴⁷R⁴⁸)₂,
- 5) C≡C R⁴⁶,
- 6) (CRⁱR^j)₂OR⁴⁶,
- 7) (CRⁱR^j)₂N(R⁴⁶R⁴⁷),
- 8) (CRⁱR^j)₂C(O)R⁴⁶,
- 9) (CRⁱR^j)₂C(O)OR⁴⁶,
- 10) (CRⁱR^j)₂R⁴⁶,
- 11) (CRⁱR^j)₂S(O)O-2R⁶¹,
- 12) (CRⁱR^j)₂S(O)O-2N(R⁴⁶R⁴⁷),
- 13) OS(O)O-2R⁶¹,
- 14) N(R⁴⁶)C(O)R⁴⁷,
- 15) N(R⁴⁶)S(O)O-2R⁶¹,
- 16) (CRⁱR^j)₂N(R⁴⁶)R⁶¹,
- 17) (CRⁱR^j)₂N(R⁴⁶)R⁶¹OR⁴⁷,
- 18) (CRⁱR^j)₂N(R⁴⁶)(CR^kR^l)₂C(O)N(R⁴⁷R⁴⁸),
- 19) N(R⁴⁶)(CRⁱR^j)₂R⁶¹,
- 20) N(R⁴⁶)(CRⁱR^j)₂N(R⁴⁷R⁴⁸),
- 21) (CRⁱR^j)₂C(O)N(R⁴⁷R⁴⁸), or
- 22) oxo, or

b) a heteroaryl ring selected from the group consisting of

- a 5-membered unsaturated monocyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting of N, O or S,
- a 6-membered unsaturated monocyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting of N, O and S, and
- a 9- or 10-membered unsaturated bicyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting of N, O or S;

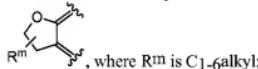
wherein any stable S heteroaryl ring atom is unsubstituted or mono- or di-substituted with oxo, and any stable C or N heteroaryl ring atom is independently unsubstituted or substituted with

- 1) halogen,
- 2) NO₂,
- 3) CN,
- 4) CR⁴⁶=C(R⁴⁷R⁴⁸)₂,
- 5) C≡CR⁴⁶,
- 6) (CRⁱR^j)_tOR⁴⁶,
- 7) (CRⁱR^j)_tN(R⁴⁶R⁴⁷),
- 8) (CRⁱR^j)_tC(O)R⁴⁶,
- 9) (CRⁱR^j)_tC(O)OR⁴⁶,
- 10) (CRⁱR^j)_tR⁴⁶,
- 11) (CRⁱR^j)_tS(O)O-2R⁶¹,
- 12) (CRⁱR^j)_tS(O)O-2N(R⁴⁶R⁴⁷),
- 13) OS(O)O-2R⁶¹,
- 14) N(R⁴⁶)C(O)R⁴⁷,
- 15) N(R⁴⁶)S(O)O-2R⁶¹,
- 16) (CRⁱR^j)_tN(R⁴⁶)R⁶¹,
- 17) (CRⁱR^j)_tN(R⁴⁶)R⁶¹OR⁴⁷,
- 18) (CRⁱR^j)_tN(R⁴⁶)(CR^kR^l)_sC(O)N(R⁴⁷R⁴⁸),
- 19) N(R⁴⁶)(CRⁱR^j)_tR⁶¹,
- 20) N(R⁴⁶)(CRⁱR^j)_tN(R⁴⁷R⁴⁸),
- 21) (CRⁱR^j)_tC(O)N(R⁴⁷R⁴⁸), or
- 22) oxo;

R², R⁸, R⁹ and R¹⁰ are independently selected from:

- 1) hydrogen,
- 2) halogen,

3) NO₂,
4) CN,
5) CR43=C(R44R45),
6) C≡CR43,
7) (CRERF₁)_nOR43,
8) (CRERF₁)_nN(R43R44),
9) (CRERF₁)_nC(O)R43,
10) (CRERF₁)_nC(O)OR43,
11) (CRERF₁)_nR43,
12) (CRERF₁)_nS(O)0-2R60,
13) (CRERF₁)_nS(O)0-2N(R43R44),
14) OS(O)0-2R60,
15) N(R43)C(O)R44,
16) N(R43)S(O)0-2R60,
17) (CRERF₁)_nN(R43)R60,
18) (CRERF₁)_nN(R43)R60OR44,
19) (CRERF₁)_nN(R43)(CRgrh)_qC(O)N(R44R45),
20) N(R43)(CRERF₁)_nR60,
21) N(R43)(CRERF₁)_nN(R44R45), and
22) (CRERF₁)_nC(O)N(R43R44),
or R2 and R8 are independently as defined above, and R9 and R10, together with the atoms to which they are attached, form the ring



, where R^m is C₁₋₆alkyl;

R¹ is selected from the group consisting of

- 1) hydrogen,
- 2) (CR^aRb)_nR40
- 3) (CR^aRb)_nOR40,
- 4) (CR^aRb)_nN(R40R41),
- 5) (CR^aRb)_nN(R40)C(O)OR41,
- 6) (CR^aRb)_nN(R40)(CRErd)₂N(R41)C(O)R49,
- 7) C₃₋₈ cycloalkyl,
- 8) (CR^aRb)_nC(O)OR40,
- 9) (CR^aRb)_nN(R40)(CRErd)₁₋₃R41,
- 10) (CR^aRb)_nS(O)0-2R6,

- 11) $(CRArB)_nS(O)_{0-2}N(R^{40}R^{41})$,
- 12) $(CRArB)_nN(R^{40})R^6OR^{41}$,
- 13) $(CRArB)_nN(R^{40})(CRCRD)_{0-6}C(O)N(R^{41}R^{42})$;

or R^1 is absent when z is a double bond

R^5 is selected from the group consisting of

- 1) C 1-6 alkyl,
- 2) =O
- 3) aryl
- 4) C₃-10 cycloalkyl
- 5) C₁₋₆alkylene-C(O)R¹¹,
- 6) C₁₋₆alkylene-C(O)R¹³
- 7) C(O)R¹¹,
- 8) C(O)R¹³,
- 9) C(O)OR¹¹,
- 10) C(O)OR¹³,
- 11) C(O)N(R¹¹R¹¹),
- 12) C(O)N(R¹³R¹³),
- 13) C(O)N(R¹¹R¹³),
- 14) CN,
- 15) NHC(O)R¹¹,
- 16) NHC(O)CF₃, and
- 17) NHC(O)C₂₋₆alkyl,

or R^1 and R^5 , together with atoms to which they are attached, form



where t is 0, 1, 2, or 3, and R^n is selected from the group consisting of hydrogen, -OR₉, NR₉R₉, C(O)NR₉R₉, or C(O)OR₉, wherein R₉ and R₉ are independently selected from the group consisting of C₁₋₆ alkyl and aryl;

R^{11} is selected from the group consisting of

- 1) aryl, and
- 2) an unsubstituted or substituted heterocyclic ring consisting of a 4-6 membered unsaturated or saturated monocyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting N, O and S, and a 9- or 10-membered unsaturated or saturated bicyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting of N, O or S; and

R¹³ is selected from the group consisting of

- 1) C₁-₆alkyl,
- 2) C₁-₆alkyloxy,
- 3) C₁-₆alkenyl,
- 4) C₁-₆alkynyl, and
- 5) CF₃;

R^a, R^b, R^c, R^d, R^e, R^f, R^g, R^h, Rⁱ, R^j, R^k, and R^l are independently selected from the group consisting of:

- 1) hydrogen,
- 2) C₁-C₆ alkyl,
- 3) halogen,
- 4) aryl,
- 5) R⁸⁰,
- 6) C₃-C₁₀ cycloalkyl, and
- 7) OR⁴,

said alkyl, aryl, and cycloalkyl being unsubstituted, monosubstituted with R⁷, disubstituted with R⁷ and R¹⁵, trisubstituted with R⁷, R¹⁵ and R¹⁶, or tetrasubstituted with R⁷, R¹⁵, R¹⁶ and R¹⁷;

R⁴, R⁴⁰, R⁴¹, R⁴², R⁴³, R⁴⁴, R⁴⁵, R⁴⁶, R⁴⁷, R⁴⁸, R⁴⁹, R⁵¹, and R⁵² are independently selected from:

- 1) hydrogen,
- 2) C₁-C₆ alkyl,
- 3) C₃-C₁₀ cycloalkyl,
- 4) aryl,
- 5) R⁸¹,
- 6) CF₃,
- 7) C₂-C₆ alkenyl, and
- 8) C₂-C₆ alkynyl,

said alkyl, aryl, and cycloalkyl is unsubstituted, mono-substituted with R¹⁸, di-substituted with R¹⁸ and R¹⁹, tri-substituted with R¹⁸, R¹⁹ and R²⁰, or tetra-substituted with R¹⁸, R¹⁹, R²⁰ and R²¹;

R⁶, R⁶⁰, R⁶¹, and R⁶³ are independently selected from:

- 1) C₁-C₆ alkyl,
- 2) aryl,
- 3) R⁸³, and
- 4) C₃-C₁₀ cycloalkyl;

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said alkyl, aryl, and cycloalkyl is unsubstituted, mono-substituted with R26, di-substituted with R26 and R27, tri-substituted with R26, R27 and R28, or tetra-substituted with R26, R27, R28 and R29;

R7, R15, R16, R17, R18, R19, R20, R21, R26, R27, R28, and R29 are independently selected from:

- 1) C1-C6 alkyl,
- 2) halogen,
- 3) OR⁵¹,
- 4) CF₃,
- 5) aryl,
- 6) C₃-C₁₀ cycloalkyl,
- 7) R⁸⁴,
- 8) S(O)₀₋₂N(R⁵¹R⁵²),
- 9) C(O)OR⁵¹,
- 10) C(O)R⁵¹,
- 11) CN,
- 12) C(O)N(R⁵¹R⁵²),
- 13) N(R⁵¹)C(O)R⁵²,
- 14) S(O)₀₋₂R⁶³,
- 15) NO₂, and
- 16) N(R⁵¹R⁵²);

R80, R81, R83 and R84 are independently selected from a group of unsubstituted or substituted heterocyclic rings consisting of a 4-6 membered unsaturated or saturated monocyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting N, O and S, and a 9- or 10-membered unsaturated or saturated bicyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting of N, O or S; and

n, p, q, r, and s are independently 0, 1, 2, 3, 4, 5 or 6, provided that, when R⁹ is hydrogen, A is substituted as defined above.

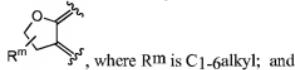
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2. (Canceled) A compound of Claim 1, or a pharmaceutically acceptable salt thereof, wherein A is an aryl ring selected from phenyl, unsubstituted or substituted as in Claim 1, or a heteroaryl ring, unsubstituted or substituted as in Claim 1, selected from the group consisting of pyridine, pyrimidine, pyrazine, pyridazine, indole, pyrrolopyridine, benzimidazole, benzoxazole, benzothiazole, and benzodiazole;

R2, R8, R9 and R10 are independently selected from the group consisting of:

- 1) hydrogen,
- 2) halogen,
- 3) OR⁴³, and
- 4) (CR^aRF₂)_pR⁴³,

or R² and R⁸ are independently as defined above, and R⁹ and R¹⁰, together with the atoms to which they are attached, form the ring



R¹ is selected from the group consisting of

- 1) hydrogen,
- 2) (CR^aRb)₁₋₂R⁴⁰
- 3) (CR^aRb)₁₋₂OR⁴⁰,
- 4) (CR^aRb)₁₋₂N(R⁴⁰R⁴¹),
- 5) (CR^aRb)₁₋₂N(R⁴⁰)C(O)OR⁴¹,
- 6) (CR^aRb)₁₋₂N(R⁴⁰)(CR^cRd)₂N(R⁴¹)C(O)R⁴⁹,
- 7) (CR^aRb)₁₋₂C(O)OR⁴⁰,
- 8) (CR^aRb)₁₋₂N(R⁴⁰)(CR^cRd)₁₋₃R⁴¹, and
- 9) cyclopropyl,

or R¹ and R⁵, together with atoms to which they are attached, form



where t is 0, 1, 2, or 3, and Rⁿ is selected from the group consisting of hydrogen, -OR_p, NR_pR_q, C(O)NR_pR_q, or C(O)OR_p, wherein RP and Rq are independently selected from the group consisting of C₁₋₆ alkyl and aryl.

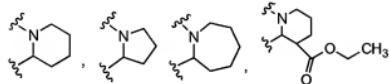
3. (Canceled) A compound of Claim 2, or a pharmaceutically acceptable salt thereof, wherein R², R⁸, R⁹, and R¹⁰ are independently selected from the group consisting of hydrogen and -OR⁴³.

4. (Canceled) A compound of Claim 3, or a pharmaceutically acceptable salt thereof, wherein A is selected from the group consisting of A is phenyl, fluorophenyl and chlorophenyl.

5. (Canceled) A compound of Claim 4, or a pharmaceutically acceptable salt thereof, wherein

R¹ is selected from the group consisting of C1-6alkyl and C3-10 cycloalkyl, or R¹ is absent when z is a double bond;

R⁵ is selected from the group consisting of C1-6 alkyl, =O, aryl, and C3-10 cycloalkyl; or R¹ and R⁵ together with the atoms to which they are attached, form



6. (Withdrawn) A compound of Claim 5, or a pharmaceutically acceptable salt thereof, selected from the group consisting of

5-(3-fluorophenyl)-3-methoxy-5,5a,6,7,8,9-hexahydro-11H-pyrido[2,1-b]quinazolin-11-one,

(5,6-cis)-5-(3-fluorophenyl)-3-methoxy-11-oxo-5,6,7,8,9,11-hexahydro-5aH-pyrido[2,1-b]quinazoline-6-carboxylate,

ethyl (5,6-cis)-11-oxo-5-phenyl-5,6,7,8,9,11-hexahydro-5aH-pyrido[2,1-b]quinazoline-6-carboxylate,

7-methoxy-2,3-dimethyl-1-phenyl-2,3-dihydroquinazolin-4(1H)-one,

6-methoxy-4-phenyl-2,3,3a,4-tetrahydropyrrolo[2,1-b]quinazolin-9(1H)-one,

3-methoxy-5-phenyl-5,5a,6,7,8,9-hexahydro-11H-pyrido[2,1-b]quinazolin-11-one,

3-methoxy-5-phenyl-5a,6,7,8,9,10-hexahydroazepino[2,1-b]quinazolin-12(5H)-one,

7-methoxy-2-methyl-4-oxo-1-phenyl-1,4-dihydroquinazolin-1-iium chloride,

2-tert-butyl-7-methoxy-1-phenylquinazolin-4(1H)-one,

2-cyclohexyl-7-methoxy-1-phenylquinazolin-4(1H)-one, and

3-cyclopropyl-7-methoxy-1-phenylquinazoline-2,4(1H,3H)-dione.

7. (Withdrawn) A method of treating a condition in a mammal, the treatment of which is effected or facilitated by Kv1.5 inhibition, which comprises administering a compound of Claim 1 in an amount that is effective at inhibiting Kv1.5.

8. (Withdrawn) A method of Claim 7, wherein the condition is cardiac arrhythmia.

9. (Withdrawn) A method of Claim 8, wherein the cardiac arrhythmia is atrial fibrillation.

10. (Withdrawn) A method of Claim 8, wherein the cardiac arrhythmia is selected from the group consisting of atrial flutter, atrial arrhythmia and supraventricular tachycardia.

11. (Withdrawn) A method of preventing a condition in a mammal, the prevention of which is effected or facilitated by Kv1.5 inhibition, which comprises administering a compound of Claim 1 in an amount that is effective at inhibiting Kv1.5.

12. (Withdrawn) A method of Claim 11, wherein the condition is cardiac arrhythmia.

13. (Withdrawn) A method of Claim 12, wherein the cardiac arrhythmia is atrial fibrillation.

14. (Withdrawn) A method of Claim 12, wherein the cardiac arrhythmia is selected from the group consisting of atrial flutter, atrial arrhythmia and supraventricular tachycardia.

15. (Withdrawn) A method of Claim 11, wherein the condition is a thromboembolic event.

16. (Withdrawn) A method of Claim 15, wherein the thromboembolic event is a stroke.

17. (Withdrawn) A method of Claim 11, wherein the condition is congestive heart failure.

18. (Canceled) A pharmaceutical formulation comprising a pharmaceutically acceptable carrier and the compound Claim 1 or a pharmaceutically acceptable crystal form or hydrate thereof.

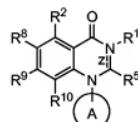
19. (Canceled) A pharmaceutical composition made by combining the compound of Claim 1 and a pharmaceutically acceptable carrier.

20. (Withdrawn) A method of treating cardiac arrhythmia comprising administering a compound of Claim 1 with a compound selected from one of the classes of compounds consisting of antiarrhythmic agents having Kv1.5 blocking activities, ACE inhibitors, angiotensin II antagonists, cardiac glycosides, L-type calcium channel blockers, T-type calcium channel blockers, selective and nonselective beta blockers, endothelin antagonists, thrombin inhibitors, aspirin, nonselective NSAIDs, warfarin, factor Xa inhibitors, low molecular weight heparin, unfractionated heparin, clopidogrel, ticlopidine, IIb/IIIa receptor antagonists, 5HT receptor antagonists, integrin receptor antagonists, thromboxane receptor antagonists, TAFI inhibitors and P2T receptor antagonists.

21. (Withdrawn) A method for inducing a condition of normal sinus rhythm in a patient having atrial fibrillation, which comprises treating the patient with a compound of Claim 1.

22. (Withdrawn) A method for treating tachycardia in a patient which comprises treating the patient with an antitachycardia device in combination with a compound of Claim 1.

23. (New) A compound of having the formula



wherein
or a pharmaceutically acceptable salt thereof, wherein
z is a single or double bond;

A is an aryl ring, wherein any stable aryl ring atom is independently unsubstituted or substituted with

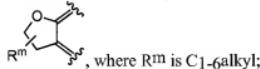
- 1) halogen,
- 2) NO₂,
- 3) CN,
- 4) CR⁴⁶=C(R⁴⁷R⁴⁸)₂,
- 5) C≡C R⁴⁶,
- 6) (CRⁱR^j)_pOR⁴⁶,
- 7) (CRⁱR^j)_pN(R⁴⁶R⁴⁷),
- 8) (CRⁱR^j)_pC(O)R⁴⁶,
- 9) (CRⁱR^j)_pC(O)OR⁴⁶,
- 10) (CRⁱR^j)_pR⁴⁶,
- 11) (CRⁱR^j)_pS(O)O-2R⁶¹,
- 12) (CRⁱR^j)_pS(O)O-2N(R⁴⁶R⁴⁷),
- 13) OS(O)O-2R⁶¹,
- 14) N(R⁴⁶)C(O)R⁴⁷,
- 15) N(R⁴⁶)S(O)O-2R⁶¹,
- 16) (CRⁱR^j)_pN(R⁴⁶)R⁶¹,
- 17) (CRⁱR^j)_pN(R⁴⁶)R⁶¹OR⁴⁷,
- 18) (CRⁱR^j)_pN(R⁴⁶)(CR^kR^l)_pC(O)N(R⁴⁷R⁴⁸),
- 19) N(R⁴⁶)(CRⁱR^j)_pR⁶¹,
- 20) N(R⁴⁶)(CRⁱR^j)_pN(R⁴⁷R⁴⁸),
- 21) (CRⁱR^j)_pC(O)N(R⁴⁷R⁴⁸), or
- 22) oxo,

R², R⁸, R⁹ and R¹⁰ are independently selected from:

- 1) hydrogen,
- 2) halogen,
- 3) NO₂,
- 4) CN,
- 5) CR⁴³=C(R⁴⁴R⁴⁵),
- 6) C≡CR⁴³,
- 7) (CR^eR^f)_pOR⁴³,
- 8) (CR^eR^f)_pN(R⁴³R⁴⁴),
- 9) (CR^eR^f)_pC(O)R⁴³,
- 10) (CR^eR^f)_pC(O)OR⁴³,

- 11) (CRE^f_jpR⁴³,
- 12) (CRE^f_jpS(O)0-2R⁶⁰,
- 13) (CRE^f_jpS(O)0-2N(R⁴³R⁴⁴),
- 14) OS(O)0-2R⁶⁰,
- 15) N(R⁴³)C(O)R⁴⁴,
- 16) N(R⁴³)S(O)0-2R⁶⁰,
- 17) (CRE^f_jpN(R⁴³)R⁶⁰,
- 18) (CRE^f_jpN(R⁴³)R⁶⁰OR⁴⁴,
- 19) (CRE^f_jpN(R⁴³)(CRgrh)_qC(O)N(R⁴⁴R⁴⁵),
- 20) N(R⁴³)(CRE^f_jpR⁶⁰,
- 21) N(R⁴³)(CRE^f_jpN(R⁴⁴R⁴⁵), and
- 22) (CRE^f_jpC(O)N(R⁴³R⁴⁴),

or R² and R⁸ are independently as defined above, and R⁹ and R¹⁰, together with the atoms to which they are attached, form the ring



R¹ is selected from the group consisting of

- 1) hydrogen,
- 2) (CRA_nRb)_nR⁴⁰
- 3) (CRA_nRb)_nOR⁴⁰,
- 4) (CRA_nRb)_nN(R⁴⁰R⁴¹),
- 5) (CRA_nRb)_nN(R⁴⁰)C(O)OR⁴¹,
- 6) (CRA_nRb)_nN(R⁴⁰)(CR^cRd)₂N(R⁴¹)C(O)R⁴⁹,
- 7) C₃-8 cycloalkyl,
- 8) (CRA_nRb)_nC(O)OR⁴⁰,
- 9) (CRA_nRb)_nN(R⁴⁰)(CR^cRd)₁₋₃R⁴¹,
- 10) (CRA_nRb)_nS(O)0-2R⁶,
- 11) (CRA_nRb)_nS(O)0-2N(R⁴⁰R⁴¹),
- 12) (CRA_nRb)_nN(R⁴⁰)R⁶OR⁴¹,
- 13) (CRA_nRb)_nN(R⁴⁰)(CR^cRd)₀₋₆C(O)N(R⁴¹R⁴²);

or R¹ is absent when z is a double bond

R⁵ is selected from the group consisting of

- 1) C 1-6 alkyl,
- 2) =O
- 3) aryl

- 4) C₃-C₁₀ cycloalkyl
- 5) C₁-C₆alkylene-C(O)R¹¹,
- 6) C₁-C₆alkylene-C(O)R¹³
- 7) C(O)R¹¹,
- 8) C(O)R¹³,
- 9) C(O)OR¹¹,
- 10) C(O)OR¹³,
- 11) C(O)N(R¹¹R¹¹),
- 12) C(O)N(R¹³R¹³),
- 13) C(O)N(R¹¹R¹³),
- 14) CN,
- 15) NHC(O)R¹¹,
- 16) NHC(O)CF₃, and
- 17) NHC(O)C₂-C₆alkyl;

R¹¹ is selected from the group consisting of

- 1) aryl, and
- 2) an unsubstituted or substituted heterocyclic ring consisting of a 4-6 membered unsaturated or saturated monocyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting N, O and S, and a 9- or 10-membered unsaturated or saturated bicyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting of N, O or S; and

R¹³ is selected from the group consisting of

- 1) C₁-C₆alkyl,
- 2) C₁-C₆alkyloxy,
- 3) C₁-C₆alkenyl,
- 4) C₁-C₆alkynyl, and
- 5) CF₃;

R^a, R^b, R^c, R^d, R^e, R^f, R^g, R^h, Rⁱ, R^j, R^k, and R^l are independently selected from the group consisting of:

- 1) hydrogen,
- 2) C₁-C₆ alkyl,
- 3) halogen,
- 4) aryl,
- 5) R⁸⁰,
- 6) C₃-C₁₀ cycloalkyl, and

7) OR⁴,

said alkyl, aryl, and cycloalkyl being unsubstituted, monosubstituted with R⁷, disubstituted with R⁷ and R¹⁵, trisubstituted with R⁷, R¹⁵ and R¹⁶, or tetrasubstituted with R⁷, R¹⁵, R¹⁶ and R¹⁷;

R⁴, R⁴⁰, R⁴¹, R⁴², R⁴³, R⁴⁴, R⁴⁵, R⁴⁶, R⁴⁷, R⁴⁸, R⁴⁹, R⁵¹, and R⁵² are independently selected from:

- 1) hydrogen,
- 2) C₁-C₆ alkyl,
- 3) C₃-C₁₀ cycloalkyl,
- 4) aryl,
- 5) R⁸¹,
- 6) CF₃,
- 7) C₂-C₆ alkenyl, and
- 8) C₂-C₆ alkynyl,

said alkyl, aryl, and cycloalkyl is unsubstituted, mono-substituted with R¹⁸, di-substituted with R¹⁸ and R¹⁹, tri-substituted with R¹⁸, R¹⁹ and R²⁰, or tetra-substituted with R¹⁸, R¹⁹, R²⁰ and R²¹;

R⁶, R⁶⁰, R⁶¹, and R⁶³ are independently selected from:

- 1) C₁-C₆ alkyl,
- 2) aryl,
- 3) R⁸³, and
- 4) C₃-C₁₀ cycloalkyl;

said alkyl, aryl, and cycloalkyl is unsubstituted, mono-substituted with R²⁶, di-substituted with R²⁶ and R²⁷, tri-substituted with R²⁶, R²⁷ and R²⁸, or tetra-substituted with R²⁶, R²⁷, R²⁸ and R²⁹;

R⁷, R¹⁵, R¹⁶, R¹⁷, R¹⁸, R¹⁹, R²⁰, R²¹, R²⁶, R²⁷, R²⁸, and R²⁹ are independently selected from:

- 1) C₁-C₆ alkyl,
- 2) halogen,
- 3) OR⁵¹,
- 4) CF₃,
- 5) aryl,
- 6) C₃-C₁₀ cycloalkyl,
- 7) R⁸⁴,
- 8) S(O)₀₋₂N(R⁵¹R⁵²),
- 9) C(O)OR⁵¹,

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- 10) C(O)R⁵¹,
- 11) CN,
- 12) C(O)N(R⁵¹R⁵²),
- 13) N(R⁵¹)C(O)R⁵²,
- 14) S(O)₀₋₂R⁶³,
- 15) NO₂, and
- 16) N(R⁵¹R⁵²);

R⁸⁰, R⁸¹, R⁸³ and R⁸⁴ are independently selected from a group of unsubstituted or

substituted heterocyclic rings consisting of a 4-6 membered unsaturated or saturated monocyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting N, O and S, and a 9- or 10-membered unsaturated or saturated bicyclic ring with 1, 2, 3 or 4 heteroatom ring atoms selected from the group consisting or N, O or S; and n, p, q, r, and s are independently 0, 1, 2, 3, 4, 5 or 6, provided that, when R⁹ is hydrogen, A is substituted as defined above;

and wherein said compound is selected from the group consisting of 7-methoxy-2,3-dimethyl-1-phenyl-2,3-dihydroquinazolin-4(1H)-one; 7-methoxy-2-methyl-4-oxo-1-phenyl-1,4-dihydroquinazolin-1-iium chloride; 2-tert-butyl-7-methoxy-1-phenylquinazolin-4(1H)-one; 2-cyclohexyl-7-methoxy-1-phenylquinazolin-4(1H)-one; and 3-Cyclopropyl-7-methoxy-1-phenylquinazoline-2,4(1H,3H)-dione.

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Claim 24 (New) A pharmaceutical formulation comprising a pharmaceutically acceptable carrier and the compound of Claim 23 or a pharmaceutically acceptable crystal form or hydrate thereof.